

IN THE UNITED STATES PATENT AND TRADEMARK OFFICE

U.S. Patent No.: 7,235,576 B1

Serial No. : 10/042,203

Filed : January 11, 2002

Issued : June 26, 2007

For : OMEGA-CARBOXYARYL SUBSTITUTED DIPHENYL UREAS AS RAF
KINASE INHIBITORS

**PETITION FOR CERTIFICATE OF CORRECTION
PURSUANT TO 37 C.F.R. § 1.322 OR § 1.323**

Commissioner for Patents
P.O. Box 1450
Alexandria, VA 22313-1450

Sir:

Applicants hereby request that the above-identified U.S. patent be corrected in accordance with the attached Certificate of Correction.

C.F.R. § 1.322

- ☒ The mistake(s) were incurred through the fault of the Patent and Trademark Office and are clearly disclosed in the records of the Office. Therefore, no fee is due.

Enclosed herewith is a Form PTO-1050 listing errors that have been found in the above-identified patent.

The issued patent does not reflect the following amendments made to the application as allowed:

- 1.) Amendments to the Specification made with the submission filed on November 19, 2002 with the Request for Continued Examination (RCE);

2.) The cancellation of claims 52, 53, 54, 69, 70, 71 and 74-109 made in the amendment filed on July 27, 2004. The Notice of Allowance dated October 16, 2006 indicates these claims are not pending. These claims appear in the patent as claims 2, 3, 4, 6, 9, 11 and 12-47.

3.) The addition of claims 122, 123, 124, 126, 129 and 131-137 in the amendments dated May 12, 2005 and October 26, 2005. These claims were entered and identified as pending and allowed in the Notice dated October 16, 2006.

In presenting this Certificate of Correction, Applicants have assumed the allowed claims (claims 68 and 122-127) would be numbered consecutively from 1-17. Therefore, claims 2-47 have been replaced with the allowed claims renumbered as claims 2-17.

Accordingly, patentees and their assignee respectfully request that the Patent and Trademark Office issue a Certificate of Correction pursuant to 37 C.F.R. § 1.322 or § 1.323, respectively.

Respectfully submitted,

/Richard J. Traverso/

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Attorney for Applicants

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Attorney Docket No.: BAYER-0025-A

Date: July 23, 2007

UNITED STATES PATENT AND TRADEMARK OFFICE

CERTIFICATE OF CORRECTION

PATENT NO : 7,235,576 B1

DATED : June 26, 2007

INVENTOR(S) : Riedl et al.

It is certified that error appears in the above-identified patent and that said Letters Patent are hereby corrected as shown below:

In column 3, line 31, reads "(CH₂)_m-and", should read -- (CH₂)_m-, and --.

In column 4, line 37, reads "methanesulphonic, should read -- methanesulfonic --.

In column 9, line 27, reads "ie.," should read -- i.e., --.

In column 9, line 39, reads "enclosure", should read -- disclosure --.

In column 17, line 1, reads "under stream", should read -- under a stream --.

In column 18, lines 39 and 40, reads "Aromatic", should read -- Aromatic --.

In column 20, line 29, reads "Methylcarbamoly)", should read -- Methylcarbamoyl --.

In column 21, line 27, reads "(N-methylcarbamoly)", should read -- (N-Methylcarbamoyl) --.

In column 21, line 38, reads "(N-methylcarbamoly)", should read -- (N-Methylcarbamoyl) --.

In column 23, line 62, reads "then mixture", should read -- then the mixture --.

In column 30, line 4, reads "(4-(4-Methylsulfonylphenyloxy)", should read -- (4-(4-Methylsulfonylphenoxy) --.


In column 35, line 58, reads "heat", should read -- heated --.

In column 36, line 44, reads "ω-(Aroylamino)phenyl", should read -- ω-(Arylamino)phenyl --.

In column 42, line 29, reads "(triflouromethyl)", should read -- trifluoromethyl --.

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Arlington, Va. 22201

PATENT NO. 7,235,576 B1



Burden Hour Statement: This form is estimated to take 1.0 hour to complete. Time will vary depending upon the needs of the individual case. Any comments on the amount of time you are required to complete this form should be sent to the Chief Information Officer, U.S. Patent and Trademark Office, Washington, DC 20231. DO NOT SEND FEES OR COMPLETED FORMS TO THIS ADDRESS. SEND TO: Assistant Commissioner for Patents, Washington, DC 20231.

UNITED STATES PATENT AND TRADEMARK OFFICE

CERTIFICATE OF CORRECTION

PATENT NO.: 7,235,576 B1

DATED: June 26, 2007

INVENTOR(S): Riedl et al.

It is certified that error appears in the above-identified patent and that said Letters Patent are hereby corrected as shown below:

In columns 92-98, replace claims 2-47 with the claims 2-17 as follows:

2. A pharmaceutically acceptable salt of claim 1, which is

a) a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid; or

b) an acid salt of an organic or inorganic base containing an alkali metal cation, an alkaline earth metal cation, an ammonium cation, an aliphatic substituted ammonium cation or an aromatic substituted ammonium cation.

3. A pharmaceutically acceptable salt of claim 1 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

4. A pharmaceutically acceptable salt which is the tosylate salt of

N-(5-*tert*-butyl-2-methoxy phenyl)-*N'*-(4-(4-methoxy-3-(*N*-methylcarbamoyl) phenoxy) phenyl) urea,

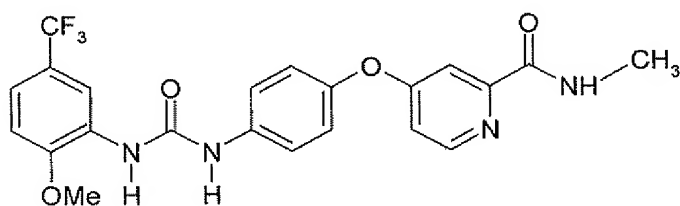
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea,

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea,

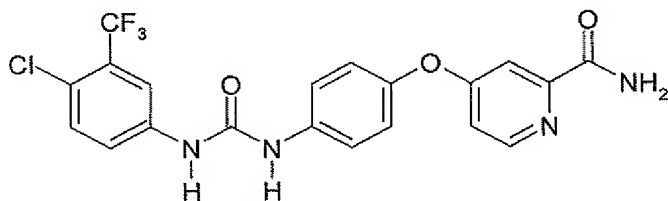
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea.

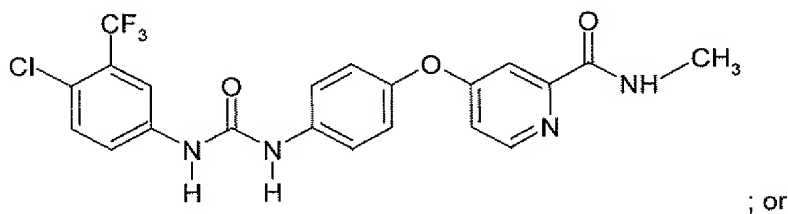
5. A pharmaceutically acceptable salt of a compound which is:
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



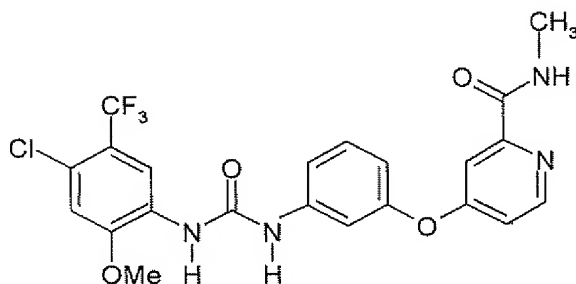
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



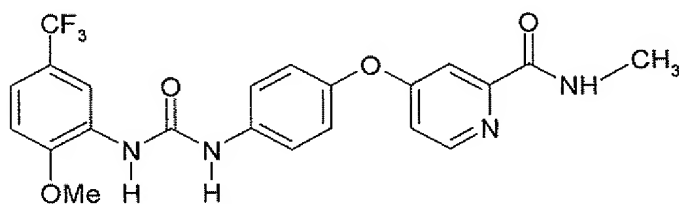
N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



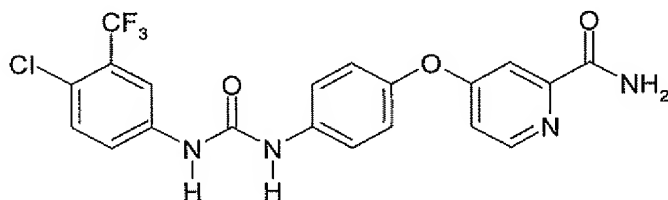
6. A pharmaceutically acceptable salt of claim 5 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

7. A pharmaceutically acceptable salt which is the tosylate salt of

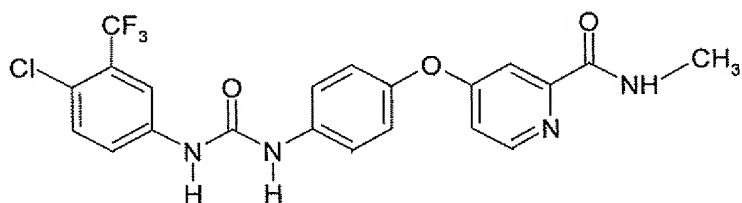
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

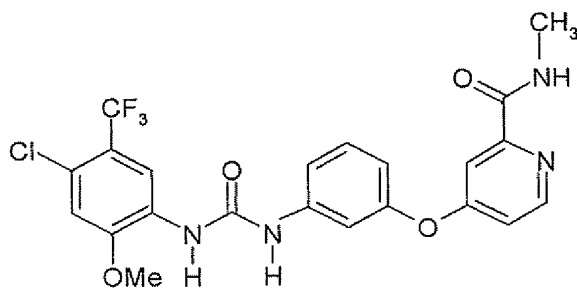


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



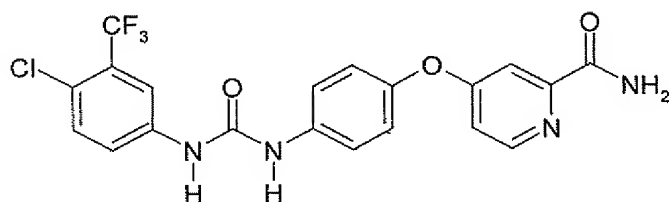
; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

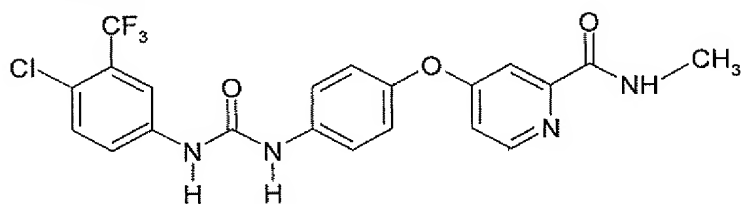


8. A pharmaceutically acceptable salt of a compound which is:

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

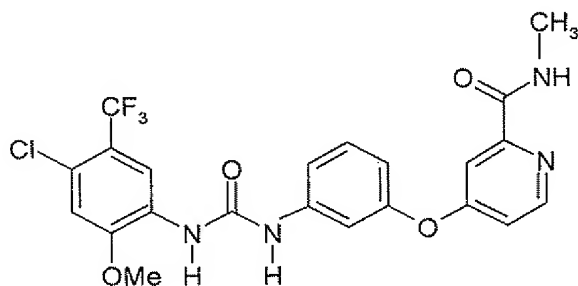


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



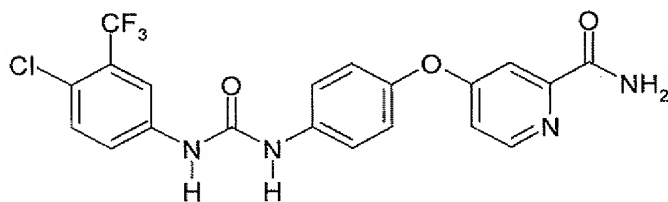
; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

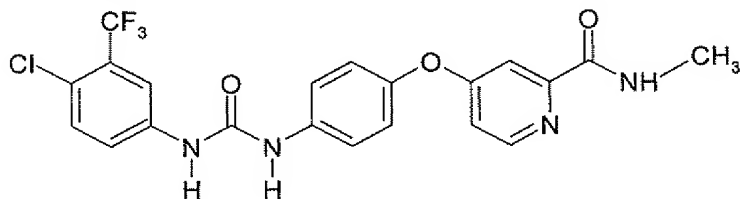


9. A pharmaceutically acceptable salt of claim 8 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

10. A pharmaceutically acceptable salt which is the tosylate salt of
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

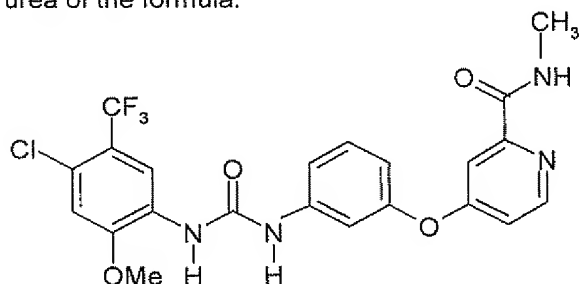


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



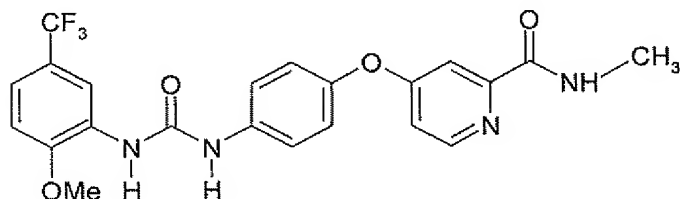
; or

N-(2-methoxy-4-chloro-5-(trifluoromethyl) phenyl)-*N'*-(3-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

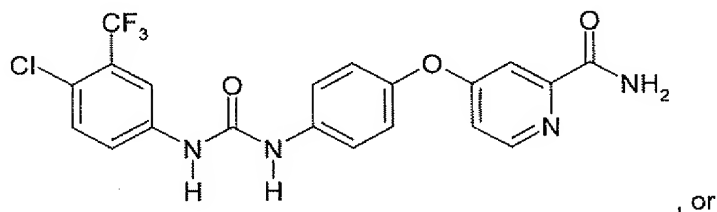


11. A pharmaceutically acceptable salt of a compound which is:

N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:

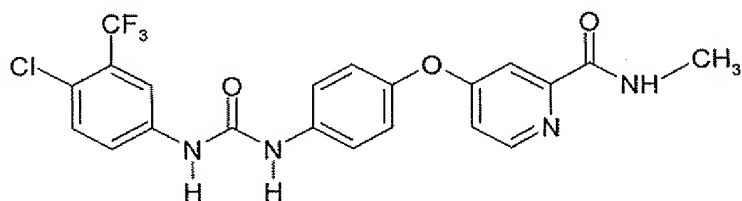


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



, or

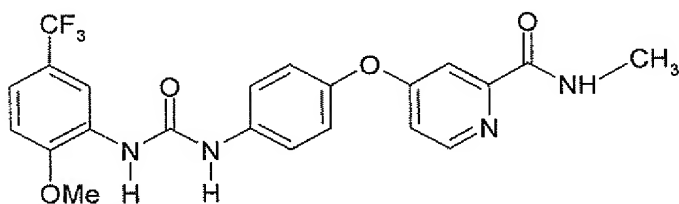
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



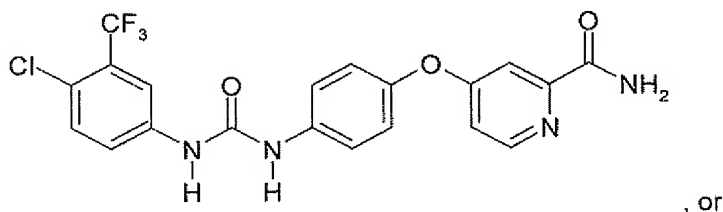
12. A pharmaceutically acceptable salt of claim 11 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

13. A pharmaceutically acceptable salt which is the tosylate salt of

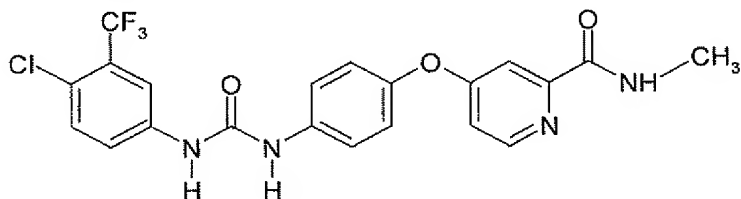
N-(2-methoxy-5-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

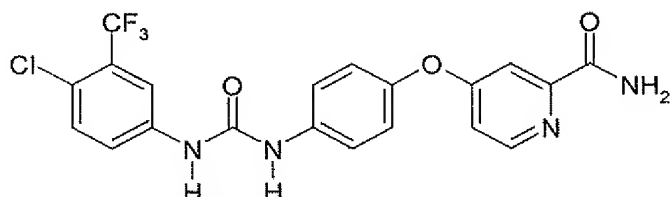


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula

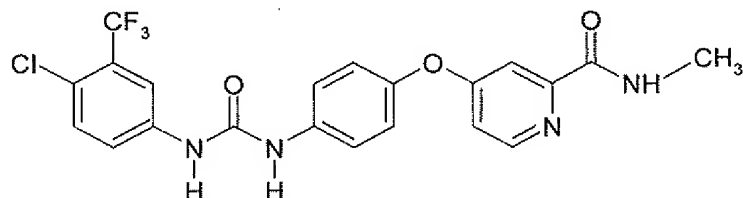


14. A pharmaceutically acceptable salt of a compound which is:

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:

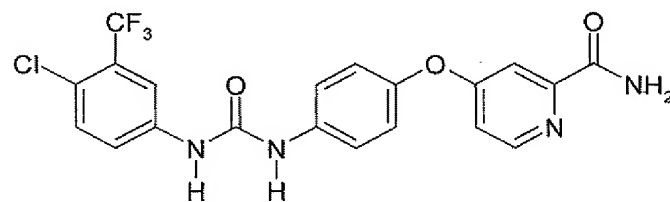


N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula:



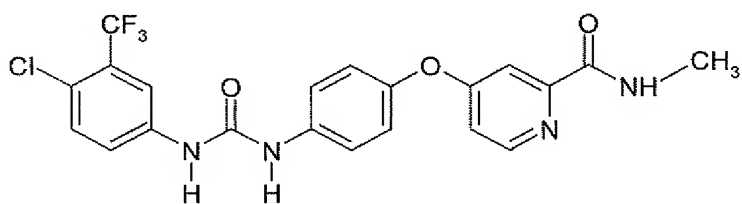
15. A pharmaceutically acceptable salt of claim 14 which is a basic salt of an organic acid or an inorganic acid which is hydrochloric acid, hydrobromic acid, sulfuric acid, phosphoric acid, methanesulfonic acid, trifluoromethanesulfonic acid, benzenesulfonic acid, p-toluene sulfonic acid (tosylate salt), 1-naphthalene sulfonic acid, 2-naphthalene sulfonic acid, acetic acid, trifluoroacetic acid, malic acid, tartaric acid, citric acid, lactic acid, oxalic acid, succinic acid, fumaric acid, maleic acid, benzoic acid, salicylic acid, phenylacetic acid, or mandelic acid.

16. A pharmaceutically acceptable salt which is the tosylate salt of
N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-carbamoyl-4-pyridyloxy) phenyl) urea of the formula:



or

N-(4-chloro-3-(trifluoromethyl) phenyl)-*N'*-(4-(2-(*N*-methylcarbamoyl)-4-pyridyloxy) phenyl) urea of the formula



- 17.** A pharmaceutical composition comprising a pharmaceutically acceptable salt of claim 1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15 or 16 and pharmaceutically acceptable carrier.